

Propagating uncertainty using FE advanced Monte-Carlo methods: application to non-linear hyperelastic models

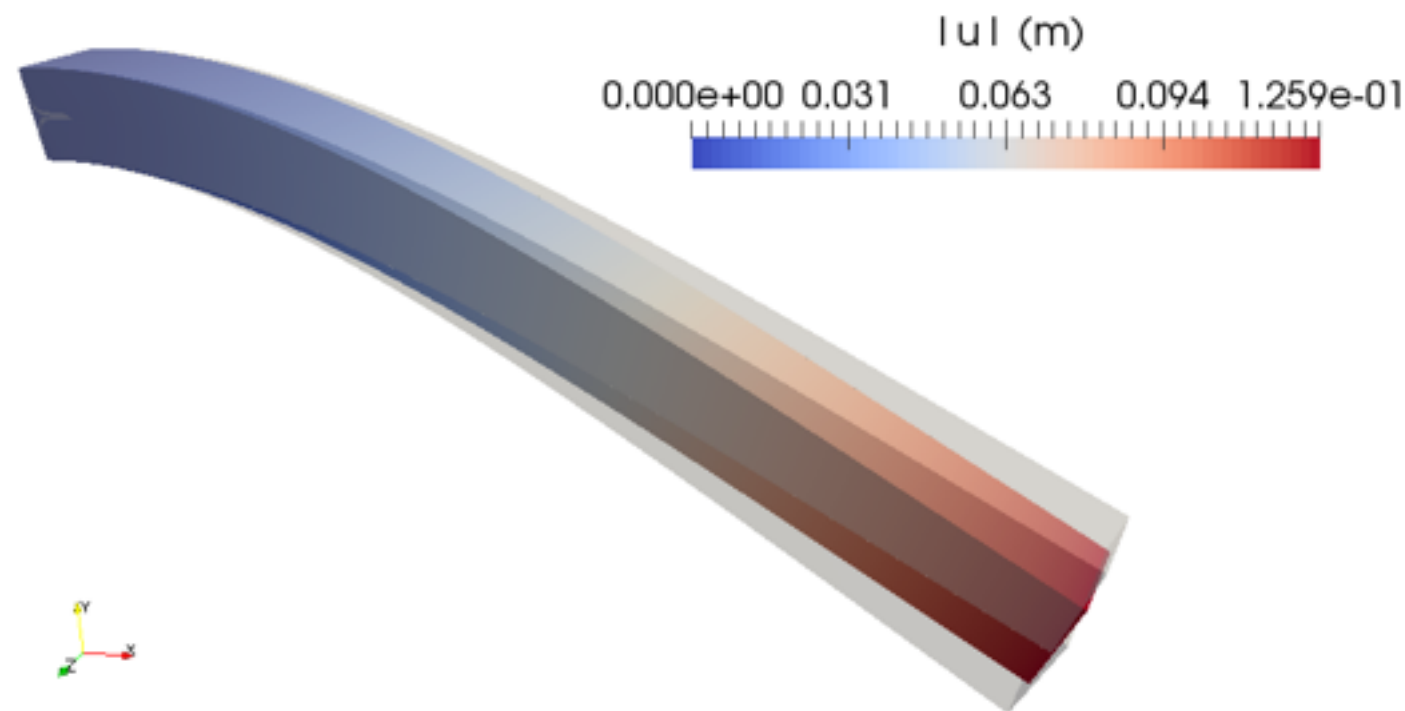
Paul Hauseux, Jack S. Hale and Stéphane P. A. Bordas

05/09/2016



Context: Soft-tissue biomechanics simulations with uncertainty

- Non-linear hyperelastic model as a stochastic PDE with **random coefficients**
- *Partially-intrusive* **Monte-Carlo methods** to propagate uncertainty



Deformation of the beam: mean +/- standard deviation

- Implementation: DOLFIN [Logg et al. 2012] and chaospy [Feinberg and Langtangen 2015]
- Ipyparallel and mpi4py to massively parallelise individual forward model runs across a cluster

1) Monte-Carlo method

- A non-linear stochastic system to solve can be written as:

$$F(\boldsymbol{u}, \boldsymbol{\omega}) = \mathbf{0}$$

- Expected value of a quantity of interest [Caflisch 1998]:

$$E(\psi(\boldsymbol{u}(\boldsymbol{x}, \boldsymbol{\omega}))) = \int_{\Omega} \psi(\boldsymbol{u}(\boldsymbol{x}, \boldsymbol{\omega})) dP(\boldsymbol{\omega}) = \frac{1}{Z} \sum_{z=1}^Z \psi(\boldsymbol{u}(\boldsymbol{x}, \boldsymbol{\omega}_z)) + o\left(\frac{\|\psi\|}{\sqrt{Z}}\right)$$

Probability space: (Ω, \mathcal{F}, P)

Random parameters: $\boldsymbol{\omega} = (\omega_1, \omega_2, \dots, \omega_M)$

- The classical Monte-Carlo approach:

$$E(\psi(\boldsymbol{u}(\boldsymbol{x}, \boldsymbol{\omega})))^{MC} \approx \frac{1}{Z} \sum_{z=1}^Z \psi(\boldsymbol{u}(\boldsymbol{x}, \boldsymbol{\omega}_z))$$

2) MC method with use of sensitivity information

- Expected value of a quantity of interest [Cao *et al.* 2004]:

$$E(\psi(\mathbf{u}(\mathbf{x}, \boldsymbol{\omega})))^{SD-MC} \approx \frac{1}{Z} \sum_{z=1}^Z \left(\psi(\mathbf{u}(\mathbf{x}, \boldsymbol{\omega}_z)) - \sum_{i=1}^M \frac{d\psi}{d\omega_i}(\bar{\boldsymbol{\omega}}) \times (\omega_i - \bar{\omega}_i) \right)$$

- Tangent linear model to evaluate the sensitivity derivatives [Farrell *et al.* 2013]:

$$\underbrace{\frac{\partial F(\mathbf{u}, \boldsymbol{\omega})}{\partial \mathbf{u}}}_{U \times U} \underbrace{\frac{d\mathbf{u}}{d\boldsymbol{\omega}}}_{U \times M} = - \underbrace{\frac{\partial F(\mathbf{u}, \boldsymbol{\omega})}{\partial \boldsymbol{\omega}}}_{U \times M}$$

U: size of the deterministic problem
M: number of random parameters

- First and Second moments of the displacement:

$$\bar{\mathbf{u}} \approx \frac{1}{Z} \sum_{z=1}^Z \left(\mathbf{u}(\mathbf{x}, \boldsymbol{\omega}_z) - \sum_{i=1}^M \frac{d\mathbf{u}}{d\omega_i}(\bar{\boldsymbol{\omega}}) \times (\omega_i - \bar{\omega}_i) \right)$$

$$\bar{\mathbf{u}}^2 \approx \frac{1}{Z} \sum_{z=1}^Z \left(\mathbf{u}^2(\mathbf{x}, \boldsymbol{\omega}_z) - 2\bar{\mathbf{u}} \sum_{i=1}^M \frac{d\mathbf{u}}{d\omega_i}(\bar{\boldsymbol{\omega}}) \times (\omega_i - \bar{\omega}_i) \right)$$

3) Multi-level MC method with use of PCE

- Polynomial chaos expansion (PCE) [Wiener 1936]:

$$u^k(x, \omega) = \sum_{\alpha \in \mathcal{J}_{m,p}} u^k_{\alpha}(x) H_{\alpha}(\omega)$$

- ML-MC method [Matthies 2008, Giles 2015]:

Algorithm 1 Algorithm for the multilevel Polynomial Chaos Expansion Monte-Carlo method

```
1: Solve the deterministic system with average parameters to obtain  $u^d$ 
2:  $k \leftarrow 1$ 
3: while no convergence do
4:   for  $z = 1$  to  $Z$  do
5:     Generate  $\omega_z = (\omega_1^z, \omega_2^z, \dots, \omega_M^z)$ 
6:     Generate  $u^k(\omega_z) = F_{pce}(u^{k-1}(\omega_z))$  or  $u^d$  if  $k == 1$ 
7:     Call to deterministic solver to do  $d$  (1 or more) iterations with starting values  $u^k(\omega_z)$ 
       and all random parameter function of  $\omega_z$ 
8:     output:  $u^k(\omega_z)$  after  $d$  iterations
9:   end for
10:  Calculate  $F_{pce}$ , the PCE of  $u^k$  from  $Z$  values of  $\omega_z$  and  $u^k(\omega_z)$ 
11:   $k = k + 1$ 
12: end while
```

4) 3D Numerical simulations

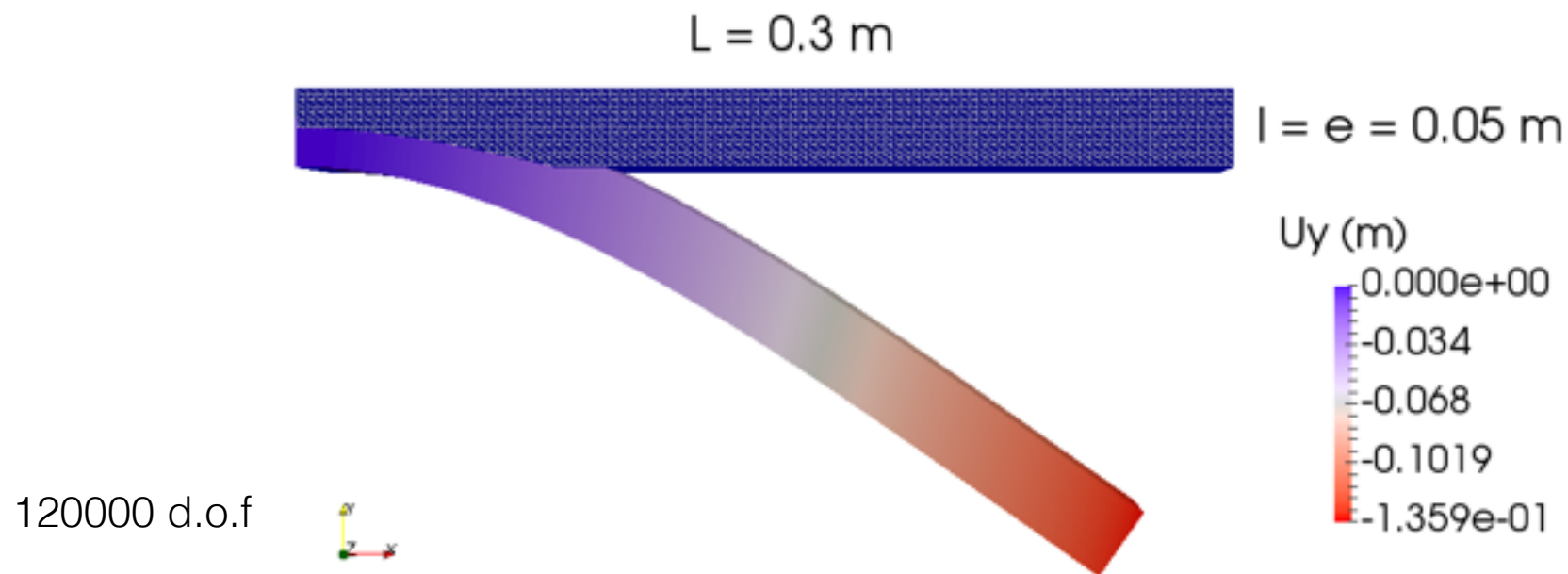


Fig: Mesh, initial configuration and deformed configuration.

- The stored strain energy density function for a compressible Mooney–Rivlin material:

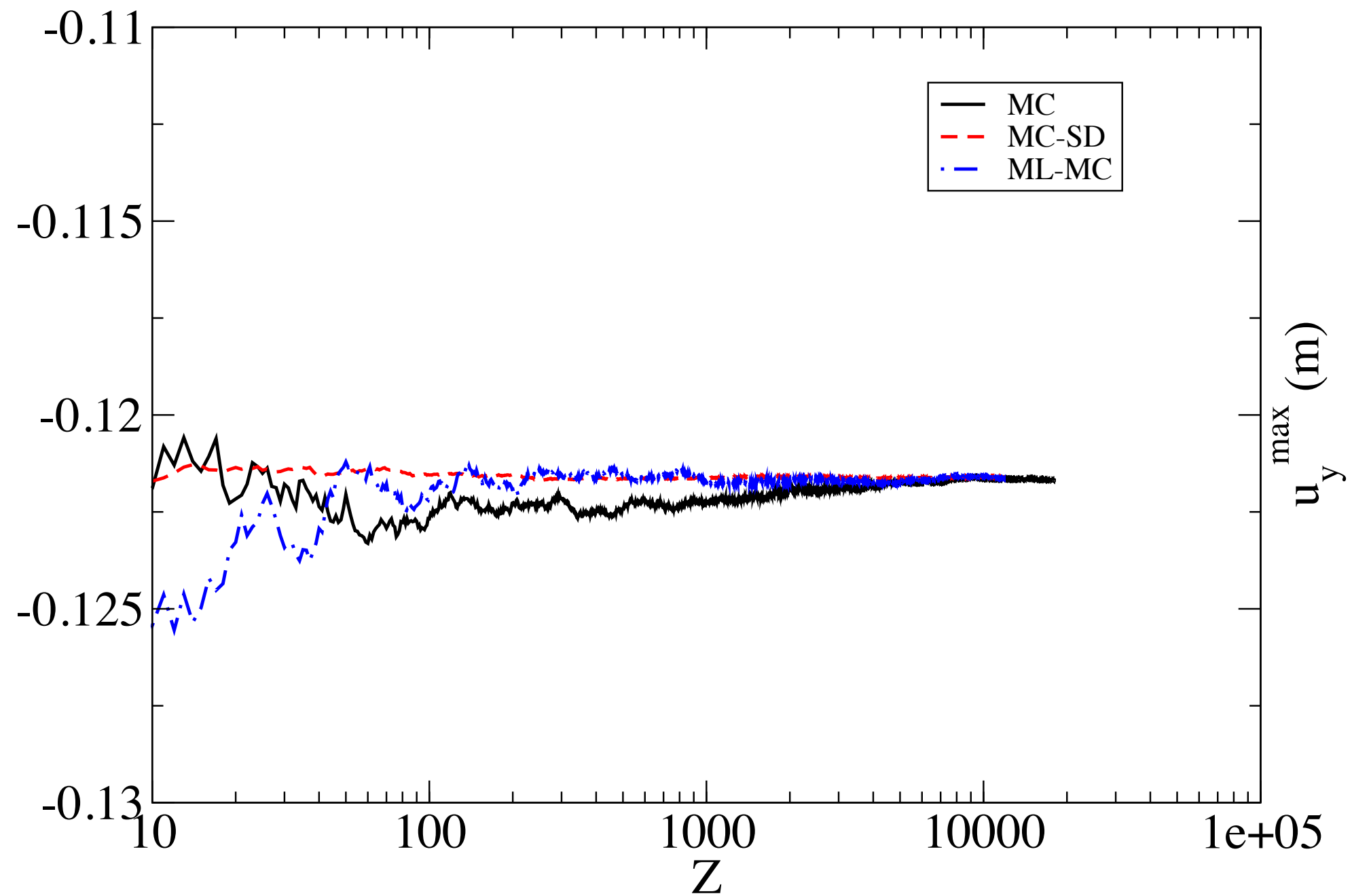
$$W = C_1(\bar{I}_1 - 3) + C_2(\bar{I}_2 - 3) + D_1(\det \mathbf{F} - 1)^2$$

- The total potential energy: $\Pi = W d\mathbf{x} - \rho \mathbf{g} d\mathbf{x}$, ($\mathbf{g} = g\vec{y}$, $g = 9.81 \text{ m.s}^{-2}$)

- 2 RV with beta(2,2) distribution:

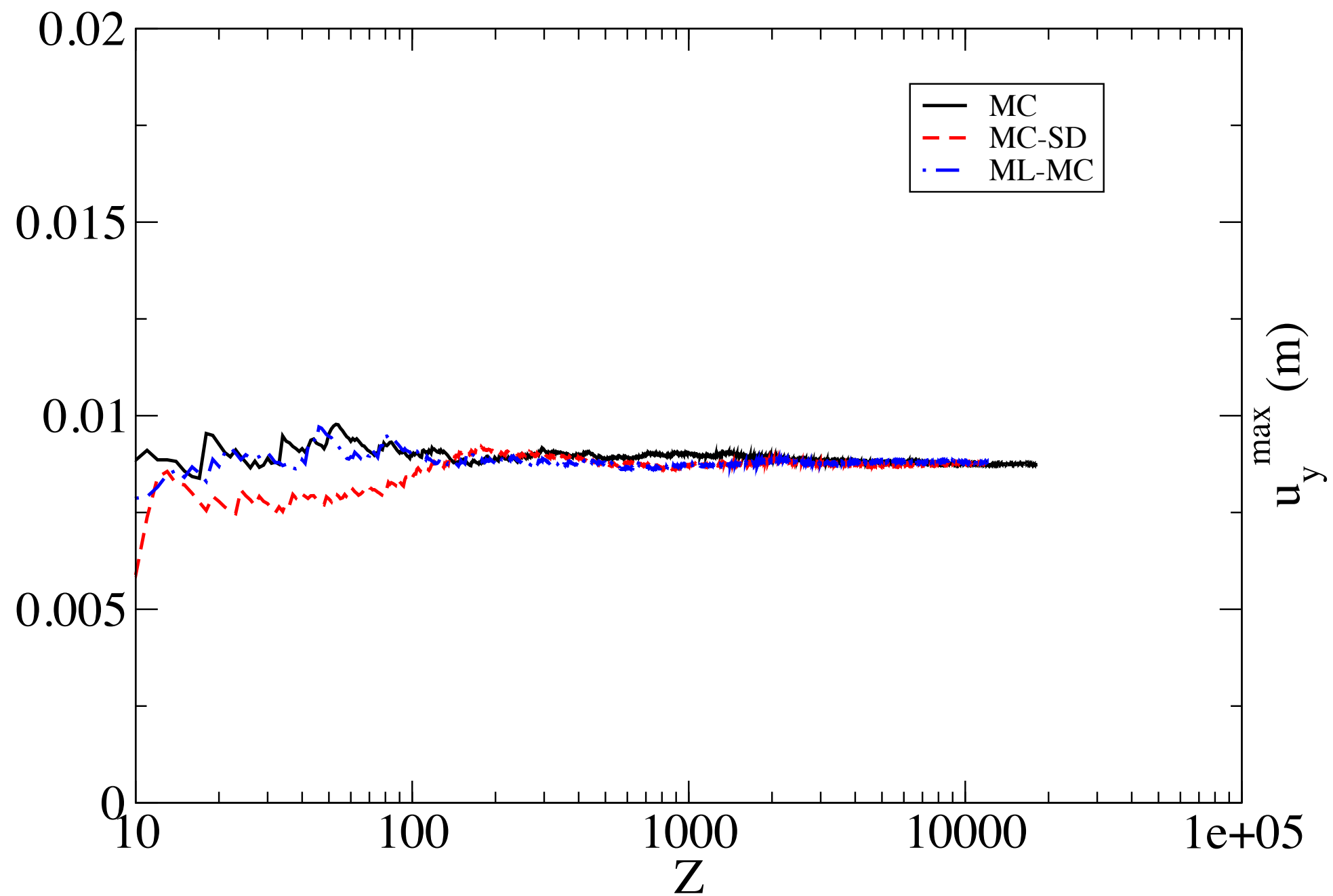
$$\begin{aligned} \rho(\omega_1) &= \rho^0(1 + \omega_1/2) \\ D_1(\omega_2) &= D_1^0(1 + \omega_2) \end{aligned} \quad \begin{cases} D_1^0 = 2 \cdot 10^5 \text{ Pa} \\ C_2 = 2 \cdot 10^5 \text{ Pa} \\ C_1 = 10^4 \text{ Pa} \\ \rho^0 = 600 \text{ kg/m}^3 \end{cases}$$

4) 3D Numerical simulations



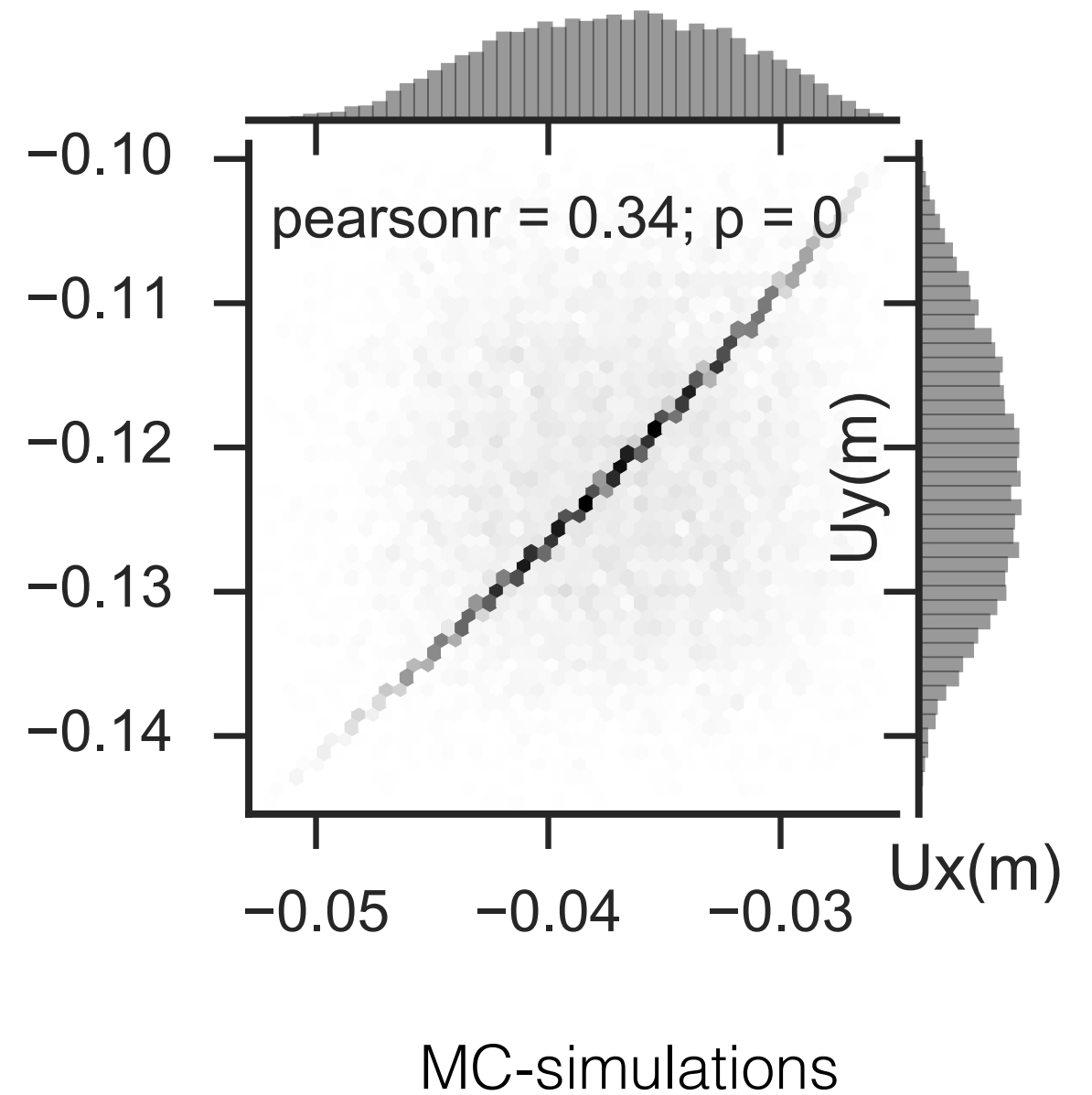
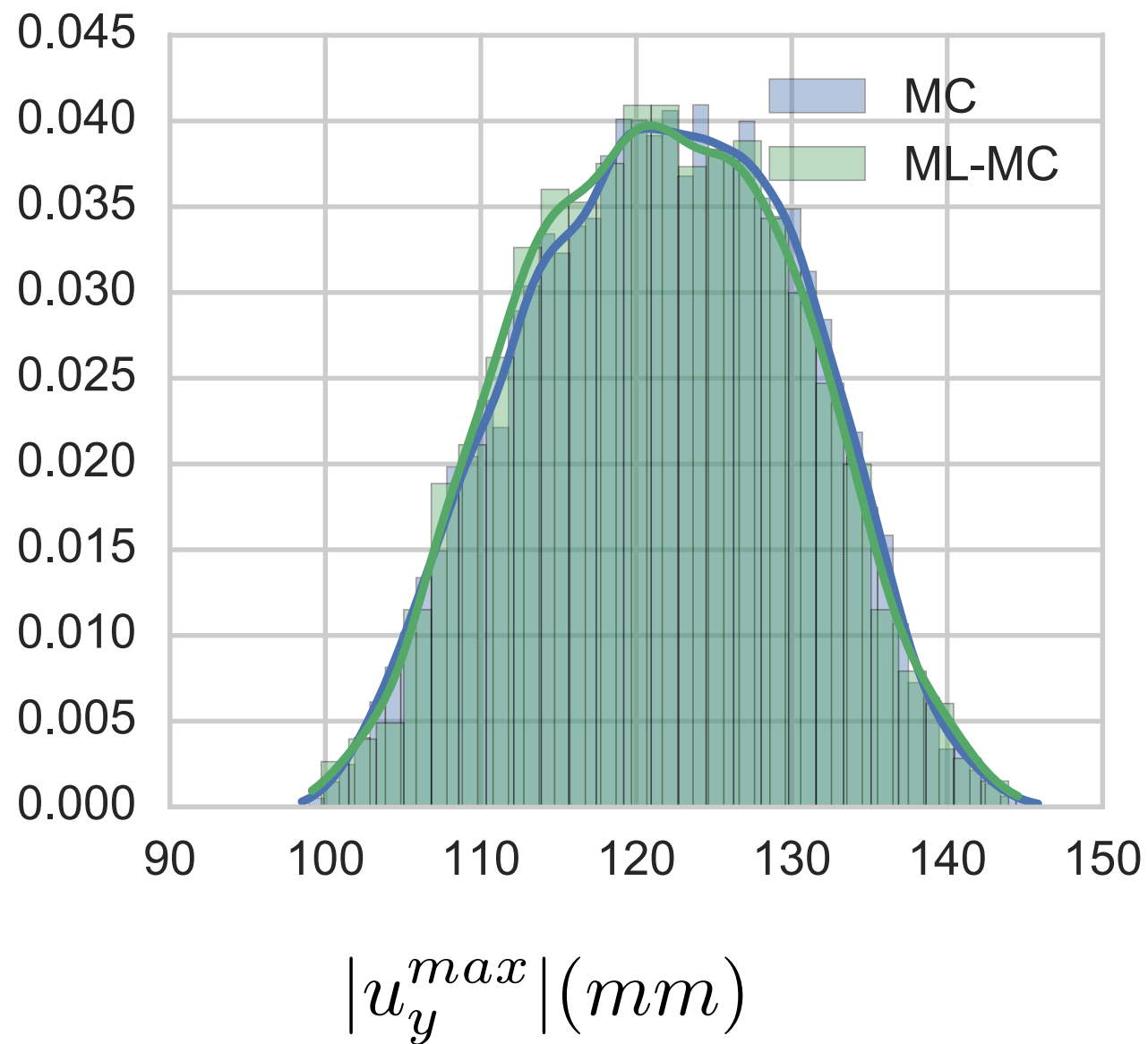
Mean

4) 3D Numerical simulations



Std

4) 3D Numerical simulations



Conclusion

- *Partially-intrusive Monte-Carlo methods* to propagate uncertainty
- By using sensitivity information and multi-level methods with polynomial chaos expansion we demonstrate that computational **workload can be reduced** by one order of magnitude over commonly used schemes
- Implementation: DOLFIN [Logg et al. 2012] and chaospy [Feinberg and Langtangen 2015]
- Ipyparallel and mpi4py to massively **parallelise** individual forward model runs across a cluster